

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092616\
 Data File : BM007574.D
 Acq On : 26 Sep 2016 16:53
 Operator : UM/SJ
 Sample : H4855-09
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H9008

Quant Time: Sep 26 17:52:10 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM092316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 26 03:32:16 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	236500	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	1120645	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.41	164	895621	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.16	188	2075133	20.00	ng/ul	0.00
75) Chrysene-d12	21.34	240	2755841	20.00	ng/ul	0.00
83) Perylene-d12	23.60	264	2499785	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.31	96	6783	1.41	ng/uL	0.00
5) Phenol-d5	6.95	99	128300	6.11	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.12	67	334789	29.66	ng/ul	0.00
9) 2-Chlorophenol-d4	7.32	132	356709	23.40	ng/ul	0.00
13) 4-Methylphenol-d8	8.48	113	259211	14.47	ng/ul	0.00
19) Nitrobenzene-d5	8.93	128	246483	30.77	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	280739	30.84	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.19	165	482810	26.62	ng/ul	0.00
29) 4-Chloroaniline-d4	10.72	131	59588	3.13	ng/ul	0.01
43) Dimethylphthalate-d6	13.82	166	2036752	31.26	ng/ul	0.00
46) Acenaphthylene-d8	14.10	160	2320043	30.69	ng/ul	0.00
51) 4-Nitrophenol-d4	14.63	143	63346	5.51	ng/ul	0.00
57) Fluorene-d10	15.40	176	1800573	31.51	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.53	200	375543	29.66	ng/ul	0.00
70) Anthracene-d10	17.25	188	2990531	31.36	ng/ul	0.00
76) Pyrene-d10	19.54	212	3615452	33.06	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.46	264	3455255	31.24	ng/ul	0.00

Target Compounds	Ovalue
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

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